

# Grid Generation and Adaptation Using an $L_p$ form of Monge-Kantorovich Optimization

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**G**rid generation and adaptation by Monge-Kantorovich optimization [1] produces optimal grids that equidistribute the local grid error, thereby minimizing the total error [2]. Optimality [1] was defined as minimizing the  $L_2$  norm

$$\|\mathbf{x}' - \mathbf{x}\|_2 = \left[ \frac{1}{2} \int |\mathbf{x}' - \mathbf{x}|^2 \rho(\mathbf{x}) d\mathbf{x} \right]^{1/2}$$

with the local constraint that the map  $\mathbf{x}' = \mathbf{x}'(\mathbf{x})$  has Jacobian  $\det[\partial(\mathbf{x}')/\partial(\mathbf{x})]$  equal to  $\rho(\mathbf{x})/\rho'(\mathbf{x}')$ , where  $\rho(\mathbf{x})$  and  $\rho'(\mathbf{x}')$  are the local errors at successive time steps (initial and target densities). In [1] it was shown that this method approximately minimizes the distortion of the map (the trace of the covariant metric tensor), and that the resulting equation, the Monge-Ampère equation, can be effectively solved by multigrid preconditioned Newton-Krylov methods.

In [3] we extended this approach to more general domains in 2D and to 3D, showing that the advantages detailed in [1] apply to these more general cases.

*Fig. 1. Target density.*

We have recently investigated the  $L_p$  form of this error equidistribution problem. That is, we minimize

$$\|\mathbf{x}' - \mathbf{x}\|_p = \left[ \frac{1}{p} \int |\mathbf{x}' - \mathbf{x}|^p \rho(\mathbf{x}) d\mathbf{x} \right]^{1/p}$$

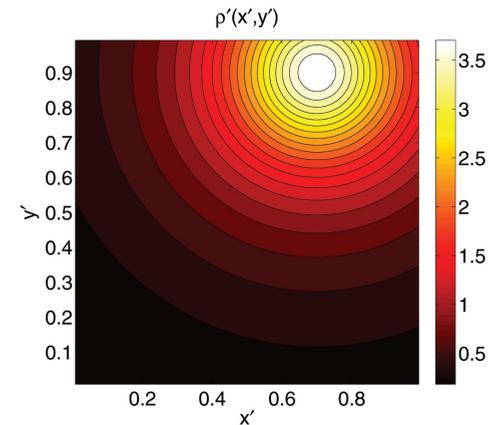
with the same Jacobian (equidistribution) constraint  $\det[\partial(\mathbf{x}')/\partial(\mathbf{x})] = \rho(\mathbf{x})/\rho'(\mathbf{x}')$ , for  $p > 1$ . The motivation for considering  $p \neq 2$  is that, relative to  $\|\mathbf{x}' - \mathbf{x}\|_p$  for  $1 < p < 2$ ,  $\|\mathbf{x}' - \mathbf{x}\|_2$  penalizes very little for small excursions but greatly for large excursions. Also, there has been a considerable amount of work in the mathematics

literature for  $p = 1$ , involving the use of the deformation method, discussed in [1]. Minimizing  $\|\mathbf{x}' - \mathbf{x}\|_p$  with the Jacobian constraint, we conclude that

$$\mathbf{x}' = \mathbf{x} + |\nabla\mu|^{\frac{2-p}{p-1}} \nabla\mu(\mathbf{x}) \tag{1}$$

where  $\mu(\mathbf{x})$  is the Legendre transform of a Lagrange multiplier enforcing the Jacobian constraint. The form corresponding to (1) for  $p = 2$  is  $\mathbf{x}' = \mathbf{x} + \nabla\mu(\mathbf{x})$ . In the difficult limit  $p \rightarrow 1$ , we find that  $\mu(\mathbf{x})$  satisfies the Eikonal equation  $|\nabla\mu| = 1$ . Substituting equation (1) into the Jacobian condition, we obtain a generalized Monge-Ampère equation for  $\mu(\mathbf{x})$ . We have found that this equation can also be solved by Newton-Krylov methods.

In Fig. 1 we show a target density  $\rho'(\mathbf{x}')$ , which is peaked at  $x = 0.7, y = 0.9$ , and the ratio of the maximum to the minimum density is  $\sim 22$ . The initial density is  $\rho(\mathbf{x}) = 1$ . In Fig. 2 we show the grid  $\mathbf{x}'$  obtained with these densities for a uniform initial grid  $\mathbf{x}$ , using various values for  $p$ . For  $p \gtrsim 1.5$ , the grids seem reasonable, concentrated near the peak of  $\rho'$  and with little distortion, i.e., the cells are close to square.



For  $p = 2$ , the grid lines are orthogonal to the boundary. For  $p < 1.5$ , and especially for  $p \rightarrow 1$ , some of the cells are seriously distorted and stretch from the origin to the peak of  $\rho'$  at  $x = 0.7, y = 0.9$ . (The darker grid lines have  $x = 0.25$ ,

$\downarrow p_c \backslash p_m \rightarrow$	1	1.01	1.02	1.05	1.1	1.25	1.5	1.75	2	2.25	2.5	Distortion
1.01	0.0000	0	0.0000	0.0004	0.0018	0.0103	0.0320	0.0572	0.0833	0.1093	0.1348	0.2501
1.02	0.0001	0.0000	0	0.0002	0.0013	0.0091	0.0295	0.0534	0.0782	0.1029	0.1270	0.2367
1.05	0.0006	0.0004	0.0002	0	0.0005	0.0063	0.0236	0.0444	0.0661	0.0878	0.1089	0.2130
1.1	0.0021	0.0017	0.0013	0.0005	0	0.0031	0.0162	0.0330	0.0510	0.0690	0.0867	0.1687
1.25	0.0092	0.0084	0.0076	0.0054	0.0028	0	0.0043	0.0132	0.0238	0.0351	0.0464	0.0716
1.5	0.0223	0.0210	0.0198	0.0164	0.0118	0.0036	0	0.0020	0.0063	0.0118	0.0178	0.0158
1.75	0.0334	0.0319	0.0304	0.0264	0.0208	0.0096	0.0017	0	0.0011	0.0035	0.0068	0.0018
2	0.0425	0.0408	0.0392	0.0348	0.0285	0.0155	0.0050	0.0010	0	0.0006	0.0022	0.0014
2.25	0.0499	0.0482	0.0464	0.0417	0.0349	0.0207	0.0085	0.0029	0.0006	0	0.0004	0.0069
2.5	0.0561	0.0542	0.0525	0.0475	0.0404	0.0254	0.0118	0.0052	0.0019	0.0004	0	0.0154

Table 1. Values of the  $p_m$  norm and the mean distortion for calculations with  $p = p_c$

$x = 0.75, y = 0.25, y = 0.75$  and are shown to emphasize the distortion for small  $p$ .) Further, the grid lines are far from orthogonal to the boundary. This adverse behavior near the boundary and especially near the corner  $(0, 0)$  for  $p \rightarrow 1$  is traced to the fact that for  $p = 1, \mu$  satisfies the Eikonal equation, which leads to a boundary layer for  $p$  close to unity.

More quantitative measures of the grid quality are contained in Table 1. The rows are labeled by  $p_c$ , which specifies the value of  $p$  used in the computations. The columns are labeled by  $p_m$ , the value of  $p$  for which the norm is measured. For example, for the highlighted in red entry with  $p_c = 1.01$  and  $p_m = 2$ , we computed the map solving the generalized Monge-Ampère equation with  $p = p_c$ , but measured  $\|x' - x\|_p$  for  $p = p_m$ . (The norms are scaled to the diagonal value with  $p_m = p_c$  with unity subtracted.) Note that in all cases, the terms below the diagonal are smaller than the transposed values  $p_m \leftrightarrow p_c$ . That is, computing the grid with large  $p_c$  works well even when quality is measured by small  $p_m$ , whereas computing with small  $p_c$  leads to large norms  $\|x' - x\|_{p_m}$  for large  $p_m$ . It appears that  $p_c = 2$  works well, but from this measure larger  $p_c$  seems slightly better. Note, however, that the measured distortion (mean of the trace of the metric tensor) is minimal at  $p_c = 2$ , orders of magnitude larger for small  $p_c$  and a factor of 10 larger for  $p = 2.5$ . Overall, these results indicate that  $p_c = 2$  leads to the highest quality grids.

Based on these results, we conclude that grid generation and adaptation works best with  $p = 2$ . We have also observed that the domain of attraction of Newton’s method is largest for  $p = 2$  and becomes much smaller for  $p$  close to unity. Further, there is a well-developed preconditioning strategy for  $p = 2$  [1].

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 [2] G. Lapenta, *J. Comput. Phys.* **193**, 159 (2004).  
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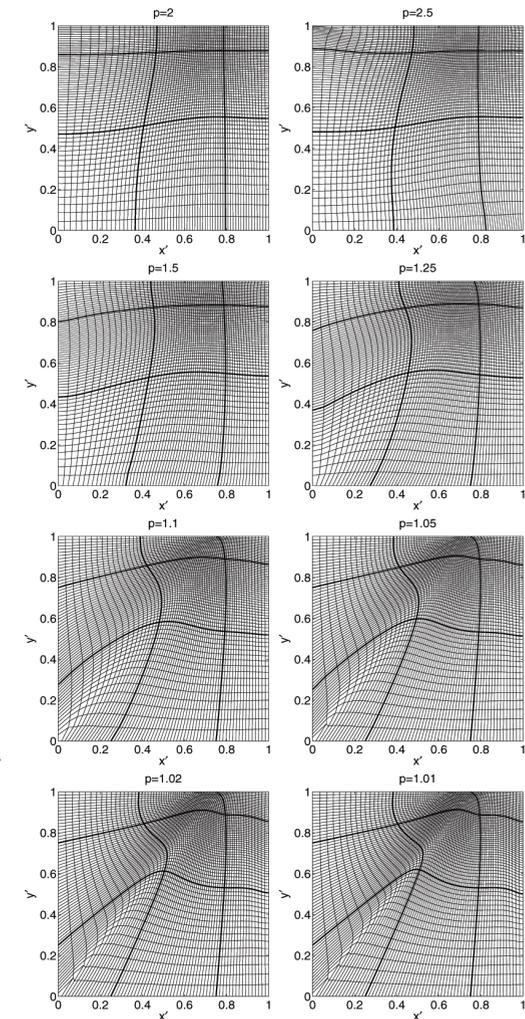


Fig. 2. Grids for various  $p$ .

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